AMENDMENT

Subject matter to be added is in bold and underlined.

Subject matter to be deleted is in bold and strikethrough.

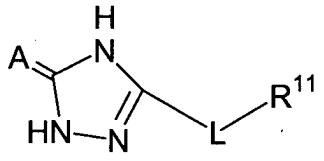
In the Claims:

Please cancel claims 11-12 without prejudice or disclaimer.

Please enter new claims 19-25 as follows.

This listing of claims will replace all prior versions and listings of claims in the application.

1. (Original) A compound of formula (I):



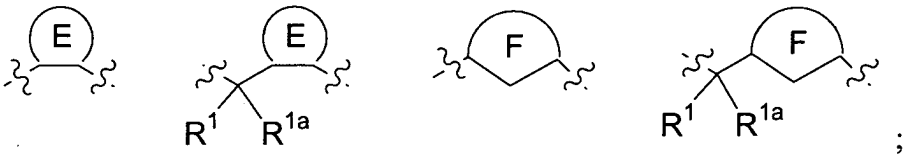
(I)

or a stereoisomer or pharmaceutically acceptable salt form thereof, wherein;

A is O or S;

L is $-(CR^2R^3)-(CR^4R^5)_n-(CR^6R^7)_{n1}-$;

alternatively, L is selected from the group:



;

R^1 is, independently at each occurrence, H, C_{1-6} alkyl, C_{2-6} alkenyl, or C_{2-6} alkynyl;

R^{1a} is, independently at each occurrence, H, C_{1-6} alkyl, C_{2-6} alkenyl, or C_{2-6} alkynyl;

R^2 is Q^1, C_{1-6} alkylene-Q^1, C_{2-6} alkenylene-Q^1, C_{2-6} alkynylene-Q^1, $-(CR^aR^{a1})_rO(CR^aR^{a1})_s-Q^1$, $-(CR^aR^{a1})_rNR^a(CR^aR^{a1})_s-Q^1$,

$-(CR^aR^{a1})_rC(O)(CR^aR^{a1})_s-Q^1$, $-(CR^aR^{a1})_rC(O)O(CR^aR^{a1})_s-Q^1$, $-(CR^aR^{a1})_rOC(O)(CR^aR^{a1})_s-Q^1$, $-(CR^aR^{a1})_rC(O)NR^aR^{a1}$, $-(CR^aR^{a1})_rC(O)NR^a(CR^aR^{a1})_s-Q^1$, $-(CR^aR^{a1})_rNR^aC(O)(CR^aR^{a1})_s-Q^1$, $-(CR^aR^{a1})_rOC(O)O(CR^aR^{a1})_s-Q^1$, $-(CR^aR^{a1})_rOC(O)NR^a(CR^aR^{a1})_s-Q^1$, $-(CR^aR^{a1})_rNR^aC(O)NR^a(CR^aR^{a1})_s-Q^1$, $-(CR^aR^{a1})_rS(O)_p(CR^aR^{a1})_s-Q^1$, $-(CR^aR^{a1})_rSO_2NR^a(CR^aR^{a1})_s-Q^1$, $-(CR^aR^{a1})_rNR^aSO_2(CR^aR^{a1})_s-Q^1$, or $-(CR^aR^{a1})_rNR^aSO_2NR^a(CR^aR^{a1})_s-Q^1$;

R^3 is Q, C_{1-6} alkylene-Q, C_{2-6} alkenylene-Q, C_{2-6} alkynylene-Q, $-(CR^aR^{a1})_rO(CR^aR^{a1})_s-Q$, $-(CR^aR^{a1})_rNR^a(CR^aR^{a1})_s-Q$, $-(CR^aR^{a1})_rC(O)(CR^aR^{a1})_s-Q$, $-(CR^aR^{a1})_rC(O)O(CR^aR^{a1})_s-Q$, $-(CR^aR^{a1})_rC(O)NR^aR^{a1}$, $-(CR^aR^{a1})_rC(O)NR^a(CR^aR^{a1})_s-Q$, $-(CR^aR^{a1})_rNR^aC(O)(CR^aR^{a1})_s-Q$, $-(CR^aR^{a1})_rS(O)_p(CR^aR^{a1})_s-Q$, $-(CR^aR^{a1})_rSO_2NR^a(CR^aR^{a1})_s-Q$, or $-(CR^aR^{a1})_rNR^aSO_2(CR^aR^{a1})_s-Q$;

Q is, independently at each occurrence, H, CHF_2, CH_2F, CF_3, a C_{3-13} carbocycle substituted with 0-5 R^d, or a 5-14 membered heterocycle consisting of carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and substituted with 0-5 R^d;

Q^1 is, independently at each occurrence, H, a C_{3-13} carbocycle substituted with 0-5 R^d, or a 5-14 membered heterocycle consisting of carbon atoms and 1-4 heteroatoms selected from the group consisting of N, NR^{10}, O, and $S(O)_p$, and substituted with 0-5 R^d;

R^4 is H, C_{1-6} alkyl substituted with 0-1 R^b, C_{2-6} alkenyl substituted with 0-1 R^b, or C_{2-6} alkynyl substituted with 0-1 R^b;

R^5 is H, C_{1-6} alkyl substituted with 0-1 R^b, C_{2-6} alkenyl substituted with 0-1 R^b, or C_{2-6} alkynyl substituted with 0-1 R^b;

R^6 is H, C_{1-6} alkyl substituted with 0-1 R^b, C_{2-6} alkenyl substituted with 0-1 R^b, or C_{2-6} alkynyl substituted with 0-1 R^b;

R^7 is H, C_{1-6} alkyl substituted with 0-1 R^b, C_{2-6} alkenyl substituted with 0-1 R^b, or C_{2-6} alkynyl substituted with 0-1 R^b;

n is 0 or 1;

n1 is 0 or 1;

alternatively, R^2 and R^3, together with the carbon atom to which they are attached, combine to form a 3-8 membered carbocyclic or heterocyclic ring consisting of carbon atoms and 0-2 ring heteroatoms selected from O, N, NR^{10}, and $S(O)_p$, and 0-2 double bonds, and substituted with 0-3 R^9; and the carbocyclic or heterocyclic ring is optionally fused to a 5-6 membered carbocycle or heterocycle consisting of carbon atoms and 0-2 ring heteroatoms selected from O, N, NR^{10}, and $S(O)_p$, and 0-2 double bonds, and substituted with 0-3 R^9;

alternatively, when n is 1, R^4 and R^5, together with the carbon atom to which they are attached, combine to form a 3-8 membered carbocyclic or heterocyclic ring consisting of carbon atoms and 0-2 ring heteroatoms selected from O, N, NR^{10}, and $S(O)_p$, and 0-2 double bonds, and substituted with 0-3 R^9; and the carbocyclic or heterocyclic ring is optionally fused to a 5-6 membered carbocycle or heterocycle consisting of carbon atoms and 0-2 ring heteroatoms selected from O, N, NR^{10}, and $S(O)_p$, and 0-2 double bonds, and substituted with 0-3 R^9;

alternatively, when n1 is 1, R^6 and R^7, together with the carbon atom to which they are attached, combine to form a 3-8 membered carbocyclic or heterocyclic ring consisting of carbon atoms and 0-2 ring heteroatoms selected from O, N, NR^{10}, and $S(O)_p$, and 0-2 double bonds, and substituted with 0-3 R^9; and the carbocyclic or heterocyclic ring is optionally fused to a 5-6 membered carbocycle or heterocycle consisting of carbon atoms and 0-2 ring heteroatoms selected from O, N, NR^{10}, and $S(O)_p$, and 0-2 double bonds, and substituted with 0-3 R^9;

ring E is a 3-8 membered carbocyclic or heterocyclic ring consisting of carbon atoms and 0-2 ring heteroatoms selected from O, N, NR^{10}, and $S(O)_p$, and 0-2 double bonds, and substituted with 0-3 R^c; and the carbocyclic or heterocyclic ring is optionally fused to a 5-6 membered carbocycle or heterocycle consisting of carbon atoms and 0-2 ring heteroatoms selected from O, N, NR^{10}, and $S(O)_p$, and 0-2 double bonds, and substituted with 0-3 R^9;

ring F is a 4-8 membered carbocyclic or heterocyclic ring consisting of carbon atoms and 0-2 ring heteroatoms selected from O, N, NR^{10}, and $S(O)_p$, and 0-2 double bonds, and substituted with 0-3 R^c; and the carbocyclic or heterocyclic ring is optionally fused to a 5-6 membered carbocycle or heterocycle consisting of carbon atoms and 0-2 ring heteroatoms selected from O, N, NR^{10}, and $S(O)_p$, and 0-2 double bonds, and substituted with 0-3 R^9;

R^{11} is –W-U-X-Y-Z-U^a-X^a-Y^a-Z^a;

W is $(CR^aR^{a1})_m$, C_{2-3} alkenylene, or C_{2-3} alkynylene;

U is O, C(O), $CR^a(OH)$, C(O)O, OC(O), $C(O)NR^{a1}$, $NR^{a1}C(O)$, OC(O)O, $OC(O)NR^{a1}$, $NR^{a1}C(O)O$, $NR^{a1}C(O)NR^{a1}$, $S(O)_p$, $S(O)_pNR^{a1}$, $NR^{a1}S(O)_p$, or $NR^{a1}SO_2NR^{a1}$;

X is absent or is C_{1-3} alkylene, C_{2-3} alkenylene, or C_{2-3} alkynylene;

Y is absent or is O, NR^{a1}, $S(O)_p$, or C(O);

Z is a C_{3-13} carbocycle substituted with 0-5 R^b, or a 5-14 membered heterocycle consisting of carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$ and substituted with 0-5 R^b;

U^a is absent or is O, NR^{a1}, C(O), $CR^a(OH)$, C(O)O, OC(O), $C(O)NR^{a1}$, $NR^{a1}C(O)$, OC(O)O, $OC(O)NR^{a1}$, $NR^{a1}C(O)O$, $NR^{a1}C(O)NR^{a1}$, $S(O)_p$, $S(O)_pNR^{a1}$, $NR^{a1}S(O)_p$, or $NR^{a1}SO_2NR^{a1}$;

X^a is absent or is C_{1-10} alkylene, C_{2-10} alkenylene, or C_{2-10} alkynylene;

Y^a is absent or is O, NR^{a1}, $S(O)_p$, or C(O);

Z^a is H, a C_{3-13} carbocycle substituted with 0-5 R^c, or a 5-14 membered heterocycle consisting of carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$ and substituted with 0-5 R^c;

provided that U, Y, Z, U^a, Y^a, and Z^a do not combine to form a N-N, N-O, O-N, O-O, $S(O)_p$-O, O-$S(O)_p$ or $S(O)_p$-$S(O)_p$ group;

R^a is, independently at each occurrence, H, C_{1-6} alkyl, phenyl, or benzyl;

R^{a1} is, independently at each occurrence, H, C_{1-6} alkyl substituted with 0-1 R^{c1}, C_{2-6} alkenyl substituted with 0-1 R^{c1}, C_{2-6} alkynyl substituted with 0-1 R^{c1}, or -$(CH_2)_r$-3-8 membered carbocyclic or heterocyclic ring consisting of carbon atoms and 0-2 ring heteroatoms selected from N, NR^{a2}, O, and $S(O)_p$, and substituted with 0-3 R^{c1};

alternatively, R^a and R^{a1} when attached to a nitrogen, together with the nitrogen to which they are attached, combine to form a 5 or 6 membered heterocycle consisting of carbon atoms and from 0-1 additional heteroatoms selected from N, NR^{a2}, O, and $S(O)_p$;

R^{a2} is, independently at each occurrence, C_{1-4} alkyl, phenyl, or benzyl;

R^{a3} is, independently at each occurrence, H, C_{1-6} alkyl substituted with R^{c1}, C_{2-6} alkenyl substituted with 0-1 R^{c1}, C_{2-6} alkynyl substituted with 0-1 R^{c1}, or -$(CH_2)_r$-3-8 membered carbocyclic or heterocyclic ring consisting of carbon atoms and 0-2 ring heteroatoms selected from N, NR^{a2}, O, and $S(O)_p$, and substituted with 0-3 R^{c1};

R^b, at each occurrence, is independently selected from C_{1-6} alkyl substituted with 0-1 R^{c1}, OR^a, SR^a, Cl, F, Br, I, =O, CN, NO_2, -NR^aR^{a1}, -$C(O)R^a$, -$C(O)OR^a$, -$C(O)NR^aR^{a1}$, -$C(S)NR^aR^{a1}$, -$NR^aC(O)NR^aR^{a1}$, -$OC(O)NR^aR^{a1}$, -$NR^aC(O)OR^a$, -$S(O)_2NR^aR^{a1}$, -$NR^aS(O)_2R^{a3}$, -$NR^aS(O)_2NR^aR^{a1}$, -$OS(O)_2NR^aR^{a1}$, -$S(O)_pR^{a3}$, -CF_3, -CF_2CF_3, -CHF_2, -CH_2F, or phenyl;

R^c is, independently at each occurrence, H, OR^a, Cl, F, Br, I, =O, CN, NO_2, CF_3, -CF_2CF_3, CH_2F, CHF_2, -$(CR^aR^{a1})_rNR^aR^{a1}$, -$(CR^aR^{a1})_rC(=NCN)NR^aR^{a1}$, -$(CR^aR^{a1})_rC(=NR^a)NR^aR^{a1}$, -$(CR^aR^{a1})_rC(=NOR^a)NR^aR^{a1}$, -$(CR^aR^{a1})_rC(O)NR^aOH$,

$-(CR^aR^{a1})_rC(O)R^{a1}$, $-(CR^aR^{a1})_rC(O)OR^{a1}$, $-(CR^aR^{a1})_rC(S)OR^{a1}$, $-(CR^aR^{a1})_rC(O)NR^aR^{a1}$, $-(CR^aR^{a1})_rNR^aC(O)R^{a1}$, $-(CR^aR^{a1})_rC(S)NR^aR^{a1}$, $-(CR^aR^{a1})_rOC(O)NR^aR^{a1}$, $-(CR^aR^{a1})_rNR^aC(O)OR^{a1}$, $-(CR^aR^{a1})_rNR^aC(O)NR^aR^{a1}$, $-(CR^aR^{a1})_rS(O)_pR^{a3}$, $-(CR^aR^{a1})_rSO_2NR^aR^{a1}$, $-(CR^aR^{a1})_rNR^aSO_2R^{a3}$, $-(CR^aR^{a1})_rNR^aSO_2NR^aR^{a1}$, C_{1-6} alkyl substituted with 0-2 R^{c1}, C_{2-6} alkenyl substituted with 0-2 R^{c1}, C_{2-6} alkynyl substituted with 0-2 R^{c1}, $-(CR^aR^{a1})_r$-C_{3-10} carbocycle substituted with 0-2 R^{c1}, or $-(CR^aR^{a1})_r$-5-14 membered heterocycle consisting of carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and substituted with 0-2 R^{c1};

alternatively, when two R^c groups are attached to the same carbon atom, they form a 3-8 membered carbocyclic or heterocyclic spiro ring C substituted with 0-2 R^{c1} and consisting of carbon atoms, 0-4 ring heteroatoms selected from O, N, and $S(O)_p$, and 0-2 double bonds, provided that ring C contains other than a S-S, O-O, or S-O bond;

alternatively, when two R^c groups are attached to adjacent carbon atoms, together with the carbon atoms to which they are attached they form a 5-7 membered carbocyclic or heterocyclic ring D substituted with 0-2 R^{c1} and consisting of carbon atoms, 0-2 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and 0-3 double bonds;

R^{c1} is, independently at each occurrence, H, C_{1-4} alkyl, OR^a, Cl, F, Br, I, =O, CF_3, CN, NO_2, $-C(O)R^a$, $-C(O)OR^a$, $-C(O)NR^aR^a$, or $-S(O)_pR^a$;

R^d is, independently at each occurrence, C_{1-6} alkyl, OR^a, Cl, F, Br, I, =O, CN, NO_2, $-NR^aR^{a1}$, $-C(O)R^a$, $-C(O)OR^a$, $-C(O)NR^aR^{a1}$, $-C(S)NR^aR^{a1}$, $-NR^aC(O)NR^aR^{a1}$, $-OC(O)NR^aR^{a1}$, $-NR^aC(O)OR^a$, $-S(O)_2NR^aR^{a1}$, $-NR^aS(O)_2R^{a3}$, $-NR^aS(O)_2NR^aR^{a1}$, $-OS(O)_2NR^aR^{a1}$, $-S(O)_pR^{a3}$, $-CF_3$, $-CF_2CF_3$, C_{3-10} carbocycle, or a 5-14 membered heterocycle consisting of carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$;

R^e is, independently at each occurrence, H, C_{1-6} alkyl, C_{1-6} alkoxy, phenoxy, benzoxy, C_{3-10} carbocycle substituted with 0-2 R^{c1}, or a 5-10 membered heterocycle consisting of carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and substituted with 0-2 R^{c1};

R^9 is, independently at each occurrence, H, $-(CR^aR^{a1})_rNR^aR^{a1}$, $-(CR^aR^{a1})_rC(O)NR^aOH$, $-(CR^aR^{a1})_rC(O)(CR^aR^{a1})_sR^e$, $-(CR^aR^{a1})_rC(O)OR^{a1}$, $-(CR^aR^{a1})_rC(S)OR^{a1}$, $-(CR^aR^{a1})_rC(O)NR^aR^{a1}$, $-(CR^aR^{a1})_rNR^aC(O)R^{a1}$, $-(CR^aR^{a1})_rC(S)NR^aR^{a1}$, $-(CR^aR^{a1})_rOC(O)NR^aR^{a1}$, $-(CR^aR^{a1})_rNR^aC(O)OR^{a1}$, $-(CR^aR^{a1})_rNR^aC(O)NR^aR^{a1}$, $-(CR^aR^{a1})_rS(O)_pR^{a3}$, $-(CR^aR^{a1})_rSO_2NR^aR^{a1}$, $-(CR^aR^{a1})_rNR^aSO_2R^{a3}$, $-(CR^aR^{a1})_rNR^aSO_2NR^aR^{a1}$, C_{1-6} alkyl substituted with 0-2 R^{c1}, C_{2-6} alkenyl substituted with 0-2 R^{c1}, C_{2-6} alkynyl substituted with 0-2 R^{c1}, $-(CR^aR^{a1})_r$-C_{3-10} carbocycle substituted with 0-2 R^{c1}, or $-(CR^aR^{a1})_r$-5-10 membered heterocycle consisting of carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and substituted with 0-2 R^{c1};

R^{10} is, independently at each occurrence, H, $-(CR^aR^{a1})_tNR^aR^{a1}$, $-(CR^aR^{a1})_rC(O)NR^aOH$, $-(CR^aR^{a1})_rC(O)(CR^aR^{a1})_sR^e$, $-(CR^aR^{a1})_rC(O)OR^{a1}$, $-(CR^aR^{a1})_rC(S)OR^{a1}$, $(-CR^aR^{a1})_rC(O)NR^aR^{a1}$, $-(CR^aR^{a1})_tNR^aC(O)R^{a1}$, $-(CR^aR^{a1})_rC(S)NR^aR^{a1}$, $-(CR^aR^{a1})_tOC(O)NR^aR^{a1}$, $-(CR^aR^{a1})_tNR^aC(O)OR^{a1}$, $-(CR^aR^{a1})_tNR^aC(O)NR^aR^{a1}$, $-(CR^aR^{a1})_rS(O)_pR^{a3}$, $-(CR^aR^{a1})_rSO_2NR^aR^{a1}$, $-(CR^aR^{a1})_tNR^aSO_2R^{a3}$, $-(CR^aR^{a1})_tNR^aSO_2NR^aR^{a1}$, C_{1-6} alkyl substituted with 0-2 R^{c1}, C_{2-6} alkenyl substituted with 0-2 R^{c1}, C_{2-6} alkynyl substituted with 0-2 R^{c1}, $-(CR^aR^{a1})_r$-C_{3-10} carbocycle substituted with 0-2 R^{c1}, or $-(CR^aR^{a1})_r$-5-10 membered heterocycle consisting of carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and substituted with 0-2 R^{c1};

m, at each occurrence, is selected from 0, 1, 2 and 3;

p, at each occurrence, is selected from 0, 1, and 2;

r, at each occurrence, is selected from 0, 1, 2, 3, and 4;

s, at each occurrence, is selected from 0, 1, 2, 3, and 4; and

t, at each occurrence, is selected from 1, 2, 3, and 4;

provided that:

(i) when L is CR^2R^3 or $CR^2R^3CH_2$, Z^a is other than H;

(ii) when Z is cyclohexyl, benzodiazepinyl or a nitrogen-containing 10-membered bicyclic heteroaryl, then Z^a is other than phenyl or phenyl fused carbocycle;

(iii) when Z is phenyl, and Z^a is oxazolyl, then R^c is other than phenyl;

(iv) when Z is a C_{5-7} cycloalkyl, then R^b is other than phenyl;

(v) when A is S, and L is 1,2-phenylene, then Z^a is other than thienyl or phenyl substituted with triazolthione;

(vi) when A is S, L is CH_2, U-X-Y forms O or S, and Z is a benzopyranyl, quinazolinyl, or triazinyl ring, then Z^a is other than phenyl;

(vii) when A is S, L is 4,5,6-7-tetrahydrobenzothienyl, and U-X-Y forms C(O)NH, Z is other than 5,6,7,8-tetrahydro-benzothieno[2,3-b]pyridinyl; and

(viii) when L is 1,2-phenylene or 1,3-phenylene, then U^a-X^a-Y^a forms other than C_{1-2} alkylene or CH_2NR^{a1}.

2. (Original) A compound according to Claim 1, wherein:

R^2 is Q^1, C_{1-6} alkylene-Q^1, C_{2-6} alkenylene-Q^1, C_{2-6} alkynylene-Q^1, $-(CR^aR^{a1})_rO(CR^aR^{a1})_s$-$Q^1$, $-(CR^aR^{a1})_rNR^a(CR^aR^{a1})_s$-$Q^1$, $-(CR^aR^{a1})_rC(O)(CR^aR^{a1})_s$-$Q^1$, $-(CR^aR^{a1})_rC(O)O(CR^aR^{a1})_s$-$Q^1$, $-(CR^aR^{a1})_rOC(O)(CR^aR^{a1})_s$-$Q^1$, $-(CR^aR^{a1})_rC(O)NR^aR^{a1}$, $-(CR^aR^{a1})_rC(O)NR^a(CR^aR^{a1})_s$-$Q^1$, $-(CR^aR^{a1})_rNR^aC(O)(CR^aR^{a1})_s$-$Q^1$, $-(CR^aR^{a1})_rS(O)_p(CR^aR^{a1})_s$-$Q^1$, $-(CR^aR^{a1})_rSO_2NR^a(CR^aR^{a1})_s$-$Q^1$, or $-(CR^aR^{a1})_rNR^aSO_2(CR^aR^{a1})_s$-$Q^1$;

R^3 is Q, C_{1-6} alkylene-Q, C_{2-6} alkenylene-Q, C_{2-6} alkynylene-Q,

$-(CH_2)_rO(CH_2)_s-Q$, $-(CH_2)_rNR^a(CH_2)_s-Q$, $-(CH_2)_rC(O)(CH_2)_s-Q$, $-(CH_2)_rC(O)O(CH_2)_s-Q$, $-(CH_2)_rC(O)NR^aR^{a1}$, $-(CH_2)_rC(O)NR^a(CH_2)_s-Q$, $-(CH_2)_rNR^aC(O)(CH_2)_s-Q$, $-(CH_2)_rS(O)_p(CH_2)_s-Q$, $-(CH_2)_rSO_2NR^a(CH_2)_s-Q$, or $-(CH_2)_rNR^aSO_2(CH_2)_s-Q$;

R^4 is H, C_{1-6} alkyl substituted with 0-1 R^b, C_{2-6} alkenyl substituted with 0-1 R^b, or C_{2-6} alkynyl substituted with 0-1 R^b;

R^5 is H, C_{1-6} alkyl substituted with 0-1 R^b, C_{2-6} alkenyl substituted with 0-1 R^b, or C_{2-6} alkynyl substituted with 0-1 R^b;

R^6 is H, C_{1-6} alkyl substituted with 0-1 R^b, C_{2-6} alkenyl substituted with 0-1 R^b, or C_{2-6} alkynyl substituted with 0-1 R^b;

R^7 is H, C_{1-6} alkyl substituted with 0-1 R^b, C_{2-6} alkenyl substituted with 0-1 R^b, or C_{2-6} alkynyl substituted with 0-1 R^b;

alternatively, R^2 and R^3, together with the carbon atom to which they are attached, combine to form a 3-8 membered carbocyclic or heterocyclic ring consisting of carbon atoms and 0-2 ring heteroatoms selected from O, N, NR^{10}, and $S(O)_p$, and 0-2 double bonds, and substituted with 0-2 R^9;

alternatively, when n is 1, R^4 and R^5 together with the carbon atom to which they are attached combine to form a 3-8 membered carbocyclic or heterocyclic ring consisting of carbon atoms and 0-2 ring heteroatoms selected from O, N, NR^{10}, and $S(O)_p$, and 0-2 double bonds, and substituted with 0-2 R^9;

alternatively, when n1 is 1, R^6 and R^7 together with the carbon atom to which they are attached combine to form a 3-8 membered carbocyclic or heterocyclic ring consisting of carbon atoms and 0-2 ring heteroatoms selected from O, N, NR^{10}, and $S(O)_p$, and 0-2 double bonds, and substituted with 0-2 R^9;

ring E is a 3-8 membered carbocyclic or heterocyclic ring consisting of carbon atoms and 0-2 ring heteroatoms selected from O, N, NR^{10}, and $S(O)_p$, and 0-2 double bonds, and substituted with 0-3 R^c;

ring F is a 4-8 membered carbocyclic or heterocyclic ring consisting of carbon atoms and 0-2 ring heteroatoms selected from O, N, NR^{10}, and $S(O)_p$, and 0-2 double bonds, and substituted with 0-3 R^c;

W is $(CR^aR^{a1})_m$;

U is O, C(O), $CR^a(OH)$, C(O)O, OC(O), $C(O)NR^{a1}$, $NR^{a1}C(O)$, $S(O)_p$, $S(O)_pNR^{a1}$, or $NR^{a1}S(O)_p$;

X is absent or is C_{1-3} alkylene;

Z is a C_{3-8} cycloalkyl substituted with 0-5 R^b, a C_{3-8} cycloalkenyl substituted with 0-5 R^b, phenyl substituted with 0-5 R^b, naphthyl substituted with 0-5 R^b, or a 5-14 membered heterocycle consisting of carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$ and substituted with 0-5 R^b;

U^a is absent or is O, NR^{a1}, C(O), $CR^a(OH)$, C(O)O, $C(O)NR^{a1}$, $NR^{a1}C(O)$, $S(O)_p$, $S(O)_pNR^{a1}$, or $NR^{a1}S(O)_p$;

X^a is absent or is C_{1-4} alkylene, C_{2-4} alkenylene, or C_{2-4} alkynylene;

Y^a is absent or is O or NR^{a1};

Z^a is a C_{6-13} carbocycle substituted with 0-5 R^c, or a 5-14 membered heterocycle consisting of carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and substituted with 0-5 R^c;

R^a is, independently at each occurrence, H, C_{1-6} alkyl, phenyl, or benzyl;

R^{a1} is, independently at each occurrence, H, C_{1-6} alkyl, C_{2-6} alkenyl, C_{2-6} alkynyl, and $-(CH_2)_r$-3-8 membered carbocyclic or heterocyclic ring consisting of carbon atoms and 0-2 ring heteroatoms selected from N, NR^{a2}, O, and $S(O)_p$;

alternatively, R^a and R^{a1} when attached to a nitrogen, together with the nitrogen to which they are attached, combine to form a 5 or 6 membered heterocycle consisting of carbon atoms and from 0-1 additional heteroatoms selected from N, NR^{a2}, O, and $S(O)_p$;

R^c is, independently at each occurrence, H, OR^a, Cl, F, Br, =O, CN, NO_2, CF_3, CH_2F, CHF_2, $-CF_2CF_3$, $-(CR^aR^{a1})_rNR^aR^{a1}$, $-(CR^aR^{a1})_rC(O)R^{a1}$, $-(CR^aR^{a1})_rC(O)OR^{a1}$, $-(CR^aR^{a1})_rC(O)NR^aR^{a1}$, $-(CR^aR^{a1})_rNR^aC(O)R^{a1}$, $-(CR^aR^{a1})_rS(O)_pR^{a3}$, $-(CR^aR^{a1})_rSO_2NR^aR^{a1}$, $-(CR^aR^{a1})_rNR^aSO_2R^{a3}$, C_{1-6} alkyl substituted with 0-1 R^{c1}, C_{2-6} alkenyl substituted with 0-1 R^{c1}, C_{2-6} alkynyl substituted with 0-1 R^{c1}, $-(CH_2)_r$-C_{3-6} carbocycle substituted with 0-2 R^{c1}, or $-(CH_2)_r$-5-6 membered heterocycle consisting of carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and substituted with 0-2 R^{c1};

alternatively, when two R^c groups are attached to the same carbon atom, they form a 3-8 membered carbocyclic or heterocyclic spiro ring C substituted with 0-2 R^{c1} and consisting of carbon atoms, 0-4 ring heteroatoms selected from O, N, and $S(O)_p$, and 0-2 double bonds, provided that ring C contains other than a S-S, O-O, or S-O bond;

alternatively, when two R^c groups are attached to adjacent carbon atoms, together with the carbon atoms to which they are attached they form a 5-7 membered carbocyclic or heterocyclic ring D substituted with 0-2 R^{c1} and consisting of carbon atoms, 0-2 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and
0-3 double bonds;

R^d is, independently at each occurrence, C_{1-6} alkyl, OR^a, Cl, F, Br, =O, CN, NO_2, $-NR^aR^{a1}$, $-C(O)R^a$, $-C(O)OR^a$, $-C(O)NR^aR^{a1}$, $-S(O)_2NR^aR^{a1}$, $-NR^aS(O)_2R^{a3}$, $-S(O)_pR^{a3}$, CF_3, C_{3-6} carbocycle and a 5-6 membered heterocycle consisting of carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$;

R^9 is, independently at each occurrence, H, $-(CR^aR^{a1})_rNR^aR^{a1}$, $-(CR^aR^{a1})_rC(O)NR^aOH$, $-(CR^aR^{a1})_rC(O)(CR^aR^{a1})_sR^e$, $-(CR^aR^{a1})_rC(O)OR^{a1}$, $-(CR^aR^{a1})_rC(O)NR^aR^{a1}$, $-(CR^aR^{a1})_rNR^aC(O)R^{a1}$, $-(CR^aR^{a1})_rOC(O)NR^aR^{a1}$,

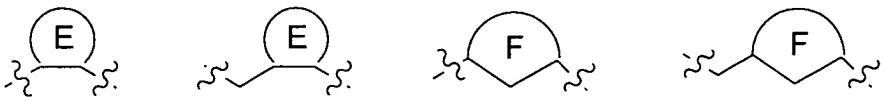
$-(CR^aR^{a1})_rNR^aC(O)OR^{a1}$, $-(CR^aR^{a1})_rS(O)_pR^{a3}$, $-(CR^aR^{a1})_rSO_2NR^aR^{a1}$, $-(CR^aR^{a1})_rNR^aSO_2R^{a3}$, C_{1-6} alkyl substituted with 0-2 R^{c1}, C_{2-6} alkenyl substituted with 0-2 R^{c1}, C_{2-6} alkynyl substituted with 0-2 R^{c1}, $-(CR^aR^{a1})_r$-C_{3-10} carbocycle substituted with 0-2 R^{c1}, or $-(CR^aR^{a1})_r$-5-10 membered heterocycle consisting of carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$ and substituted with 0-2 R^{c1}; and

R^{10} is, independently at each occurrence, H, $-(CR^aR^{a1})_tNR^aR^{a1}$, $-(CR^aR^{a1})_rC(O)NR^aOH$, $-(CR^aR^{a1})_rC(O)(CR^aR^{a1})_sR^e$, $-(CR^aR^{a1})_rC(O)OR^{a1}$, $-(CR^aR^{a1})_rC(O)NR^aR^{a1}$, $-(CR^aR^{a1})_tNR^aC(O)R^{a1}$, $-(CR^aR^{a1})_tOC(O)NR^aR^{a1}$, $-(CR^aR^{a1})_tNR^aC(O)OR^{a1}$, $-(CR^aR^{a1})_rS(O)_pR^{a3}$, $-(CR^aR^{a1})_rSO_2NR^aR^{a1}$, $-(CR^aR^{a1})_tNR^aSO_2R^{a3}$, C_{1-6} alkyl substituted with 0-2 R^{c1}, C_{2-6} alkenyl substituted with 0-2 R^{c1}, C_{2-6} alkynyl substituted with 0-2 R^{c1}, $-(CR^aR^{a1})_r$-C_{3-10} carbocycle substituted with 0-2 R^{c1}, or $-(CR^aR^{a1})_r$-5-10 membered heterocycle consisting of carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$ and substituted with 0-2 R^{c1}.

3. (Original) A compound according to Claim 2, wherein:

L is $-(CR^2R^3)$-, $-(CR^2R^3)-CH_2$-, $-(CR^2R^3)-(CH_2)_2$-, $-CH_2-(CR^4R^5)$-, or $-CH_2-(CR^4R^5)-CH_2$-;

alternatively, L is selected from the group:



;

R^2 is Q^1, C_{1-6} alkylene-Q^1, C_{2-6} alkenylene-Q^1, C_{2-6} alkynylene-Q^1, $-(CH_2)_rO(CH_2)_s-Q^1$, $-(CH_2)_rNR^a(CH_2)_s-Q^1$, $-(CH_2)_rC(O)(CH_2)_s-Q^1$, $-(CH_2)_rC(O)O(CH_2)_s-Q^1$, $-(CH_2)_rC(O)NR^a(CH_2)_s-Q^1$, $-(CH_2)_rNR^aC(O)(CH_2)_s-Q^1$, $-(CH_2)_rS(O)_p(CH_2)_s-Q^1$, $-(CH_2)_rSO_2NR^a(CH_2)_s-Q^1$, or $-(CH_2)_rNR^aSO_2(CH_2)_s-Q^1$;

R^3 is Q, C_{1-6} alkylene-Q, C_{2-6} alkenylene-Q, C_{2-6} alkynylene-Q, $-(CH_2)_rO(CH_2)_s$-Q, $-(CH_2)_rNR^a(CH_2)_s$-Q, $-(CH_2)_rC(O)(CH_2)_s$-Q, $-(CH_2)_rC(O)O(CH_2)_s$-Q, $-(CH_2)_rC(O)NR^aR^{a1}$, $-(CH_2)_rC(O)NR^a(CH_2)_s$-Q, $-(CH_2)_rNR^aC(O)(CH_2)_s$-Q, $-(CH_2)_rS(O)_p(CH_2)_s$-Q, $-(CH_2)_rSO_2NR^a(CH_2)_s$-Q, or $-(CH_2)_rNR^aSO_2(CH_2)_s$-Q;

Q is, independently at each occurrence, H, a C_{3-10} carbocycle substituted with 0-3 R^d, or a 5-10 membered heterocycle consisting of carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and substituted with 0-3 R^d;

R^4 is H or C_{1-6} alkyl;

R^5 is H or C_{1-6} alkyl;

alternatively, R^2 and R^3, together with the carbon atom to which they are attached, combine to form a 3-8 membered carbocyclic or heterocyclic ring consisting of carbon atoms, 0-2 ring heteroatoms selected from O, N, NR^{10}, and $S(O)_p$, and 0-2 double bonds, and substituted with 0-2 R^9;

alternatively, R^4 and R^5, together with the carbon atom to which they are attached, combine to form a 3-8 membered carbocyclic or heterocyclic ring consisting of carbon atoms, 0-2 ring heteroatoms selected from O, N, NR^{10}, and $S(O)_p$, and 0-2 double bonds, and substituted with 0-2 R^9;

ring E is a C_{3-7} cycloalkyl substituted with 0-2 R^c, a C_{4-7} cycloalkenyl substituted with 0-2 R^c, phenyl substituted with 0-3 R^c, or a 5-7 membered heterocyclic ring consisting of carbon atoms and 0-2 ring heteroatoms selected from O, N, NR^{10}, and $S(O)_p$, and 0-2 double bonds, and substituted with 0-3 R^c;

ring F is a C_{4-7} cycloalkyl substituted with 0-2 R^c, a C_{4-7} cycloalkenyl substituted with 0-2 R^c, phenyl substituted with 0-3 R^c, or a 5-7 membered heterocyclic ring consisting of carbon atoms and 0-2 ring heteroatoms selected from O, N, NR^{10}, and $S(O)_p$, and 0-2 double bonds, and substituted with 0-3 R^c;

U is O, C(O), C(O)NR^{a1}, $NR^{a1}C(O)$, $S(O)_p$, $S(O)_pNR^{a1}$, or $NR^{a1}S(O)_p$;

X is absent, or is methylene or ethylene;

Z is a C_{4-8} cycloalkyl substituted with 0-3 R^b, a C_{4-8} cycloalkenyl substituted with 0-3 R^b, phenyl substituted with 0-4 R^b, naphthyl substituted with 0-5 R^b, or a heterocycle substituted with 0-3 R^b and selected from the group: furanyl, tetrahydrofuranyl, thiazolyl, oxazolyl, imidazolyl, isothiazolyl, isoxazolyl, 4,5-dihydro-isoxazolyl, thienyl, triazolyl, thiadiazolyl, oxadiazolyl, pyridyl, pyrimidinyl, piperazinyl, piperidinyl, pyranyl, pyrazinyl, pyrazolyl, pyridoimidazolyl, pyrrolidinyl, pyrrolyl, indolyl, indolinyl, benzimidazolyl, benzothiazinyl, benzofuranyl, benzothiophenyl, benzoxazolyl, benzthiazolyl, benztriazolyl, benzisoxazolyl, benzisothiazolyl, quinolinyl, tetrahydroquinolinyl, isoquinolinyl, tetrahydro-isoquinolinyl, indazolyl, isobenzofuranyl, isoindazolyl, isoindolinyl, isoindolyl, methylenedioxyphenyl, and quinazolinyl;

U^a is absent or is O, NR^{a1}, C(O), C(O)NR^{a1}, $NR^{a1}C(O)$, $S(O)_p$, $S(O)_pNR^{a1}$, or $NR^{a1}S(O)_p$;

R^{a3} is, independently at each occurrence, H, C_{1-6} alkyl, C_{2-6} alkenyl, or $-(CH_2)_r$-3-8 membered carbocyclic or heterocyclic ring consisting of carbon atoms and 0-2 ring heteroatoms selected from N, NR^{a2}, O, and $S(O)_p$, and substituted with 0-3 R^{c1};

R^c is, independently at each occurrence, H, OR^a, Cl, F, Br, =O, CF_3, CH_2F, CHF_2, $-(CR^aR^{a1})_rNR^aR^{a1}$, $-(CR^aR^{a1})_rC(O)R^{a1}$, $-(CR^aR^{a1})_rC(O)OR^{a1}$, $-(CR^aR^{a1})_rC(O)NR^aR^{a1}$, $-(CR^aR^{a1})_rNR^aC(O)R^{a1}$, $-(CR^aR^{a1})_rS(O)_pR^{a3}$, $-(CR^aR^{a1})_rSO_2NR^aR^{a1}$, $-(CR^aR^{a1})_rNR^aSO_2R^{a3}$, C_{1-6} alkyl, C_{2-6} alkenyl, C_{2-6} alkynyl, C_{3-6} cycloalkyl substituted with 0-2 R^{c1}, phenyl substituted with 0-2 R^{c1}, or a 5-6 membered heterocycle consisting of carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and substituted with 0-2 R^{c1};

alternatively, when two R^c groups are attached to the same carbon atom, they form a 5-7 membered carbocyclic or heterocyclic spiro ring C substituted with 0-2 R^{c1} and

consisting of carbon atoms, 0-4 ring heteroatoms selected from O, N, and $S(O)_p$, and 0-2 double bonds, provided that ring C contains other than a S-S, O-O, or S-O bond;

alternatively, when two R^c groups are attached to adjacent carbon atoms, together with the carbon atoms to which they are attached they form a 5-7 membered carbocyclic or heterocyclic ring D substituted with 0-2 R^{c1} and consisting of carbon atoms, 0-2 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and 0-3 double bonds;

R^d is, independently at each occurrence, C_{1-6} alkyl, OR^a, Cl, F, Br, =O, $-NR^aR^{a1}$, $-C(O)R^a$, $-C(O)OR^a$, $-C(O)NR^aR^{a1}$, $-S(O)_2NR^aR^{a1}$, $-NR^aS(O)_2R^{a3}$, $-S(O)_pR^{a3}$, CF_3 or phenyl;

R^9 is, independently at each occurrence, H, $-(CR^aR^{a1})_rNR^aR^{a1}$, $-(CR^aR^{a1})_rC(O)(CR^aR^{a1})_sR^e$, $-(CR^aR^{a1})_rC(O)OR^{a1}$, $-(CR^aR^{a1})_rC(O)NR^aR^{a1}$, $-(CR^aR^{a1})_rNR^aC(O)R^{a1}$, $-(CR^aR^{a1})_rS(O)_pR^{a3}$, $-(CR^aR^{a1})_rSO_2NR^aR^{a1}$, $-(CR^aR^{a1})_rNR^aSO_2R^{a3}$, C_{1-6} alkyl substituted with 0-2 R^{c1}, C_{2-6} alkenyl substituted with 0-2 R^{c1}, C_{2-6} alkynyl substituted with 0-2 R^{c1}, $-(CR^aR^{a1})_r$-C_{3-10} carbocycle substituted with 0-2 R^{c1}, or $-(CR^aR^{a1})_r$-5-10 membered heterocycle consisting of carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and substituted with 0-2 R^{c1};

R^{10} is, independently at each occurrence, H, $-(CR^aR^{a1})_tNR^aR^{a1}$, $-(CR^aR^{a1})_rC(O)(CR^aR^{a1})_sR^e$, $-(CR^aR^{a1})_rC(O)OR^{a1}$, $-(CR^aR^{a1})_rC(O)NR^aR^{a1}$, $-(CR^aR^{a1})_tNR^aC(O)R^{a1}$, $-(CR^aR^{a1})_rS(O)_pR^{a3}$, $-(CR^aR^{a1})_rSO_2NR^aR^{a1}$, $-(CR^aR^{a1})_tNR^aSO_2R^{a3}$, C_{1-6} alkyl substituted with 0-2 R^{c1}, C_{2-6} alkenyl substituted with 0-2 R^{c1}, C_{2-6} alkynyl substituted with 0-2 R^{c1}, $-(CR^aR^{a1})_r$-C_{3-10} carbocycle substituted with 0-2 R^{c1}, or $-(CR^aR^{a1})_r$-5-10 membered heterocycle consisting of carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and substituted with 0-2 R^{c1};

r, at each occurrence, is selected from 0, 1, 2, and 3;

s, at each occurrence, is selected from 0, 1, 2, and 3; and

t, at each occurrence, is selected from 1, 2, and 3.

4. (Original) A compound according to Claim 3, wherein:

Q is, independently at each occurrence, H, a C_{3-8} carbocycle substituted with 0-3 R^d, or a 5-10 membered heterocycle consisting of carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and substituted with 0-3 R^d;

Q^1 is, independently at each occurrence, H, a C_{3-10} carbocycle substituted with 0-5 R^d, or a 5-10 membered heterocycle consisting of carbon atoms and 1-4 heteroatoms selected from the group consisting of N, NR^{10}, O, and $S(O)_p$, and substituted with 0-3 R^d;

Z is phenyl substituted with 0-3 R^b, naphthyl substituted with 0-5 R^b, pyridyl substituted with 0-3 R^b, thienyl substituted with 0-2 R^b, thiazolyl substituted with 0-2 R^b, oxazolyl substituted with 0-2 R^b, isoxazolyl substituted with 0-2 R^b, or imidazolyl substituted with 0-2 R^b;

Z^a is phenyl substituted with 0-3 R^c, naphthyl substituted with 0-3 R^c, or a heterocycle substituted with 0-3 R^c and selected from the group: furanyl, tetrahydrofuranyl, thiazolyl, oxazolyl, imidazolyl, isothiazolyl, isoxazolyl, 4,5-dihydro-isoxazolyl, thienyl, triazolyl, thiadiazolyl, oxadiazolyl, pyridyl, pyrimidinyl, piperazinyl, piperidinyl, pyranyl, pyrazinyl, pyrazolyl, pyridoimidazolyl, pyrrolidinyl, pyrrolyl, indolyl, indolinyl, benzimidazolyl, benzothiazinyl, benzofuranyl, benzothiophenyl, benzoxazolyl, benzthiazolyl, benztriazolyl, benzisoxazolyl, benzisothiazolyl, quinolinyl, tetrahydroquinolinyl, isoquinolinyl, tetrahydro-isoquinolinyl, indazolyl, isobenzofuranyl, isoindazolyl, isoindolinyl, isoindolyl, methylenedioxyphenyl, quinazolinyl, 1,1-dioxido-2,3-dihydro-4H-1,4-benzothiazin-4-yl, 1,1-dioxido-3,4-dihydro-2*H*-1-benzothiopyran-4-yl, 3,4-dihydro-2*H*-chromen-4-yl, 2*H*-chromen-4-yl, and pyrazolo[1,5-a]pyridinyl;

R^a is, independently at each occurrence, H, C_{1-6} alkyl, phenyl, or benzyl;

R^{a1} is, independently at each occurrence, H, C_{1-6} alkyl, phenyl, or benzyl;

R^{a3} is, independently at each occurrence, H, C_{1-6} alkyl, phenyl, or benzyl;

R^c is, independently at each occurrence, H, OR^a, Cl, F, Br, =O, CF_3, CH_2F, CHF_2,

$-(CR^aR^{a1})_rNR^aR^{a1}$, $-(CR^aR^{a1})_rC(O)R^{a1}$, $-(CR^aR^{a1})_rC(O)OR^{a1}$, $-(CR^aR^{a1})_rC(O)NR^aR^{a1}$, $-(CR^aR^{a1})_rNR^aC(O)R^{a1}$, $-(CR^aR^{a1})_rS(O)_pR^{a3}$, $-(CR^aR^{a1})_rSO_2NR^aR^{a1}$, $-(CR^aR^{a1})_rNR^aSO_2R^{a3}$, C_{1-6} alkyl, C_{2-6} alkenyl, C_{2-6} alkynyl, phenyl substituted with 0-2 R^{c1}, or a 5-6 membered heterocycle consisting of carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$ and substituted with 0-2 R^{c1};

alternatively, when two R^c groups are attached to the same carbon atom, they form a 5-6 membered carbocyclic or heterocyclic spiro ring C substituted with 0-2 R^{c1} and consisting of carbon atoms, 0-4 ring heteroatoms selected from O, N, and $S(O)_p$, and 0-2 double bonds, provided that ring C contains other than a S-S, O-O, or S-O bond;

alternatively, when two R^c groups are attached to adjacent carbon atoms, together with the carbon atoms to which they are attached they form a 5-6 membered carbocyclic or heterocyclic ring D substituted with 0-2 R^{c1} and consisting of carbon atoms, 0-2 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and 0-3 double bonds;

R^9 is, independently at each occurrence, H, $-(CR^aR^{a1})_rNR^aR^{a1}$, $-(CR^aR^{a1})_rC(O)(CR^aR^{a1})_sR^e$, $-(CR^aR^{a1})_rC(O)OR^{a1}$, $-(CR^aR^{a1})_rC(O)NR^aR^{a1}$, $-(CR^aR^{a1})_rNR^aC(O)R^{a1}$, $-(CR^aR^{a1})_rS(O)_pR^{a3}$, $-(CR^aR^{a1})_rSO_2NR^aR^{a1}$, $-(CR^aR^{a1})_rNR^aSO_2R^{a3}$, C_{1-6} alkyl, C_{2-6} alkenyl, C_{2-6} alkynyl, $-(CR^aR^{a1})_r$-C_{3-7} carbocycle substituted with 0-2 R^{c1}, or $-(CR^aR^{a1})_r$-5-6 membered heterocycle consisting of carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and substituted with 0-2 R^{c1}; and

R^{10} is, independently at each occurrence, H, $-(CR^aR^{a1})_tNR^aR^{a1}$, $-(CR^aR^{a1})_rC(O)(CR^aR^{a1})_sR^e$, $-(CR^aR^{a1})_rC(O)OR^{a1}$, $-(CR^aR^{a1})_rC(O)NR^aR^{a1}$, $-(CR^aR^{a1})_tNR^aC(O)R^{a1}$, $-(CR^aR^{a1})_rS(O)_pR^{a3}$, $-(CR^aR^{a1})_rSO_2NR^aR^{a1}$, $-(CR^aR^{a1})_tNR^aSO_2R^{a3}$, C_{1-6} alkyl substituted with 0-2 R^{c1}, C_{2-6} alkenyl substituted with 0-2 R^{c1}, C_{2-6} alkynyl substituted with 0-2 R^{c1}, $-(CR^aR^{a1})_r$-C_{3-10} carbocycle substituted

with 0-2 R^{c1}, or -$(CR^aR^{a1})_r$-5-10 membered heterocycle consisting of carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and substituted with 0-2 R^{c1}.

5. (Original) A compound according to Claim 4, wherein:

R^2 is Q^1, C_{1-4} alkylene-Q^1, C_{2-4} alkenylene-Q^1, C_{2-4} alkynylene-Q^1, -$(CH_2)_rO(CH_2)_s$-Q^1, -$(CH_2)_rNR^a(CH_2)_s$-Q^1, -$(CH_2)_rC(O)(CH_2)_s$-Q^1, -$(CH_2)_rC(O)O(CH_2)_s$-Q^1, -$(CH_2)_rC(O)NR^a(CH_2)_s$-Q^1, -$(CH_2)_rNR^aC(O)(CH_2)_s$-Q^1, -$(CH_2)_rS(O)_p(CH_2)_s$-Q^1, -$(CH_2)_rSO_2NR^a(CH_2)_s$-Q^1, or -$(CH_2)_rNR^aSO_2(CH_2)_s$-Q^1;

R^3 is H, C_{1-4} alkyl, C_{2-4} alkenyl, or C_{2-4} alkynyl;

R^4 is H or C_{1-4} alkyl;

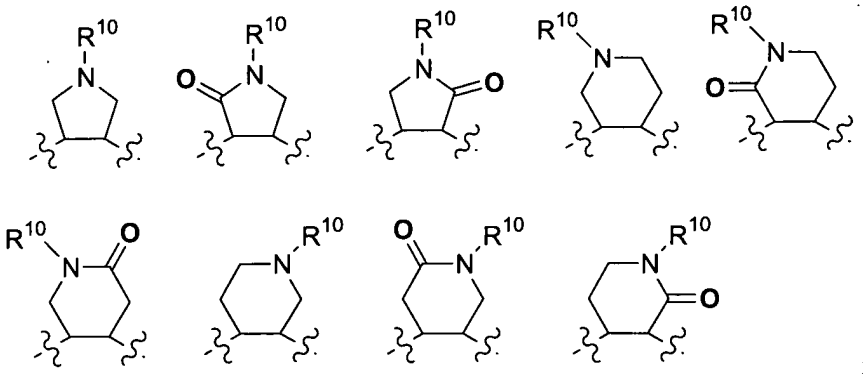
R^5 is H or C_{1-4} alkyl;

alternatively, R^2 and R^3, together with the carbon atom to which they are attached, combine to form a 3-8 membered carbocyclic or heterocyclic ring consisting of carbon atoms, 0-2 ring heteroatoms selected from O, N, NR^{10}, and $S(O)_p$, and 0-2 double bonds;

alternatively, R^4 and R^5, together with the carbon atom to which they are attached, combine to form a 3-8 membered carbocyclic or heterocyclic ring consisting of carbon atoms, 0-2 ring heteroatoms selected from O, N, NR^{10}, and $S(O)_p$, and 0-2 double bonds;

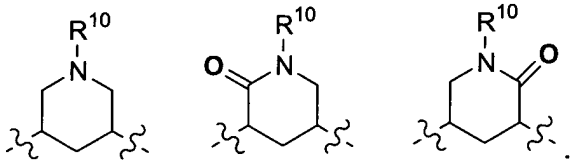
Q^1 is, independently at each occurrence, H, a C_{3-6} carbocycle substituted with 0-2 R^d, or a 5-6 membered heterocycle consisting of carbon atoms and 1-4 heteroatoms selected from the group consisting of N, NR^{10}, O, and $S(O)_p$, and substituted with 0-2 R^d;

ring E is a C_{4-7} cycloalkyl substituted with 0-2 R^c, a C_{4-7} cycloalkenyl substituted with 0-2 R^c, phenyl substituted with 0-2 R^c, or a heterocyclic ring substituted with 0-2 R^c and selected from: furanyl, tetrahydrofuranyl, thienyl, thiazolyl, oxazolyl, imidazolyl, isothiazolyl, isoxazolyl, pyrazolyl, pyrrolyl, pyridyl, pyranyl, tetrahydropyranyl, pyrimidinyl,



;

ring F a C_{4-7} cycloalkyl substituted with 0-2 R^c, a C_{4-7} cycloalkenyl substituted with 0-2 R^c, phenyl substituted with 0-2 R^c, or a heterocyclic ring substituted with 0-2 R^c and selected from: pyridyl, pyranyl, tetrahydropyranyl, pyrimidinyl,



;

U is C(O), $C(O)NR^{a1}$, $NR^{a1}C(O)$, $S(O)_p$, $S(O)_pNR^{a1}$, or $NR^{a1}S(O)_p$;

X is absent or is methylene;

Y is absent or is O;

Z is phenyl substituted with 0-3 R^b;

U^a is absent or is O;

Y^a is absent or is O;

R^a is, independently at each occurrence, H, or C_{1-4} alkyl;

R^{a1} is, independently at each occurrence, H, or C_{1-4} alkyl;

R^{a3} is, independently at each occurrence, H, C_{1-4} alkyl, phenyl, or benzyl;

R^c is, independently at each occurrence, C_{1-6} alkyl, C_{2-6} alkenyl, C_{2-6} alkynyl, OR^a, Cl, F, Br, =O, CF_3, CH_2F, CHF_2, NR^aR^{a1}, $(CR^aR^{a1})_rC(O)R^{a1}$, $(CR^aR^{a1})_rC(O)OR^{a1}$, $(CR^aR^{a1})_rC(O)NR^aR^{a1}$, $(CR^aR^{a1})_rNR^aC(O)R^{a1}$, $(CR^aR^{a1})_rS(O)_pR^{a3}$, $(CR^aR^{a1})_rSO_2NR^aR^{a1}$, $(CR^aR^{a1})_rNR^aSO_2R^{a3}$, or phenyl;

alternatively, when two R^c groups are attached to the same carbon atom, they form a 5-6 membered carbocyclic or heterocyclic spiro ring C substituted with 0-2 R^{c1} and consisting of carbon atoms, 0-4 ring heteroatoms selected from O, N, and $S(O)_p$, and 0-2 double bonds, provided that ring C contains other than a S-S, O-O, or S-O bond;

alternatively, when two R^c groups are attached to adjacent carbon atoms, together with the carbon atoms to which they are attached they form a 5-6 membered carbocyclic or heterocyclic ring D substituted with 0-1 R^{c1} and consisting of carbon atoms, 0-2 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and 0-3 double bonds;

R^e is, independently at each occurrence, H, C_{1-6} alkyl, C_{1-6} alkoxy, phenoxy, benzoxy, C_{3-6} carbocycle substituted with 0-2 R^{c1}, or a 5-6 membered heterocycle consisting of carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and substituted with 0-2 R^{c1}; and

R^{10} is, independently at each occurrence, H, $-(CH_2)_tNR^aR^{a1}$, $-(CH_2)_rC(O)(CH_2)_sR^e$, $-(CH_2)_rC(O)OR^{a1}$, $-(CH_2)_rC(O)NR^aR^{a1}$, $-(CH_2)_tNR^aC(O)R^{a1}$, $-(CH_2)_rS(O)_pR^{a3}$, $-(CH_2)_rSO_2NR^aR^{a1}$, $-(CH_2)_tNR^aSO_2R^{a3}$, C_{1-6} alkyl substituted with 0-2 R^{c1}, C_{2-6} alkenyl substituted with 0-2 R^{c1}, C_{2-6} alkynyl substituted with 0-2 R^{c1}, $-(CH_2)_r$-C_{3-10} carbocycle substituted with 0-2 R^{c1}, or $-(CH_2)_r$-5-10 membered heterocycle consisting of carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and substituted with 0-2 R^{c1}.

6. (Original) A compound according to Claim 5, wherein:

R^2 is H, C_{1-4} alkyl, C_{2-6} alkenyl, or C_{2-6} alkynyl;

R^3 is H or C_{1-4} alkylene;

R^4 is H or C_{1-4} alkyl;

R^5 is H or C_{1-4} alkyl;

alternatively, R^2 and R^3, together with the carbon atom to which they are attached, combine to form a C_{3-7} cycloalkyl, a C_{3-7} cycloalkenyl, or a 5-6 membered heterocyclic

ring consisting of carbon atoms, 1 ring heteroatom selected from O, N, NR^{10}, and $S(O)_p$, and 0-2 double bonds;

alternatively, R^4 and R^5, together with the carbon atom to which they are attached, combine to form a C_{3-7} cycloalkyl, a C_{3-7} cycloalkenyl, or a 5-6 membered heterocyclic ring consisting of carbon atoms, 1 ring heteroatom selected from O, N, NR^{10}, and $S(O)_p$, and 0-2 double bonds;

W is $(CH_2)_m$;

Y is absent;

Z is phenyl substituted with 0-1 R^b;

Z^a is phenyl substituted with 0-3 R^c, naphthyl substituted with 0-3 R^c, or a heterocycle substituted with 0-3 R^c and selected from the group: pyridyl, quinolinyl, tetrahydroquinolinyl, isoquinolinyl, tetrahydro-isoquinolinyl, imidazolyl, pyridoimidazolyl, benzimidazolyl, indolyl, indolinyl, 1,1-dioxido-2,3-dihydro-4*H*-1,4-benzothiazin-4-yl, 1,1-dioxido-3,4-dihydro-2*H*-1-benzothiopyran-4-yl, 3,4-dihydro-2*H*-chromen-4-yl, 2*H*-chromen-4-yl, pyrazolyl, and pyrazolo[1,5-a]pyridinyl;

R^b is, independently at each occurrence, C_{1-6} alkyl, OR^a, Cl, F, Br, NR^aR^{a1}, $C(O)R^a$, $C(O)OR^a$, $C(O)NR^aR^{a1}$, $S(O)_2NR^aR^{a1}$, $NR^aS(O)_2R^{a3}$, $S(O)_pR^{a3}$, or CF_3;

R^c is, independently at each occurrence, C_{1-6} alkyl, C_{2-6} alkenyl, C_{2-6} alkynyl, OR^a, Cl, F, Br, =O, CF_3, $-NR^aR^{a1}$, $-(CR^aR^{a1})_rC(O)R^{a1}$, $-(CR^aR^{a1})_rC(O)OR^{a1}$, $-(CR^aR^{a1})_rC(O)NR^aR^{a1}$, $-(CR^aR^{a1})_rNR^aC(O)R^{a1}$, $-(CR^aR^{a1})_rS(O)_pR^{a3}$, $-(CR^aR^{a1})_rSO_2NR^aR^{a1}$, or $-(CR^aR^{a1})_rNR^aSO_2R^{a3}$;

alternatively, when two R^c groups are attached to the same carbon atom, they form a 5-6 membered carbocyclic or heterocyclic spiro ring C consisting of carbon atoms, 0-4 ring heteroatoms selected from O, N, and $S(O)_p$, and 0-2 double bonds, provided that ring C contains other than a S-S, O-O, or S-O bond;

alternatively, when two R^c groups are attached to adjacent carbon atoms, together with the carbon atoms to which they are attached they form a 5-6 membered carbocyclic or

heterocyclic ring D consisting of carbon atoms, 0-2 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and 0-3 double bonds;

R^e is, independently at each occurrence, H, C_{1-6} alkyl, C_{1-6} alkoxy, phenoxy, benzoxy, phenyl substituted with 0-1 R^{c1}, or a 5-6 membered heterocycle consisting of carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and substituted with 0-1 R^{c1};

R^{10} is, independently at each occurrence, H, $-(CH_2)_rC(O)(CH_2)_sR^e$, $-(CH_2)_rC(O)OR^{a1}$, $-(CH_2)_rC(O)NR^aR^{a1}$, $-(CH_2)_rS(O)_pR^{a3}$, $-(CH_2)_rSO_2NR^aR^{a1}$, C_{1-4} alkyl substituted with 0-1 R^{c1}, C_{2-4} alkenyl substituted with 0-1 R^{c1}, C_{2-4} alkynyl substituted with 0-1 R^{c1}, $-(CH_2)_r$-C_{3-6} cycloalkyl substituted with 0-2 R^{c1}, $-(CH_2)_r$-phenyl substituted with 0-2 R^{c1}, or $-(CH_2)_r$-5-6 membered heterocycle consisting of carbon atoms and 1-4 heteroatoms selected from the group consisting of N, O, and $S(O)_p$, and substituted with 0-2 R^{c1};

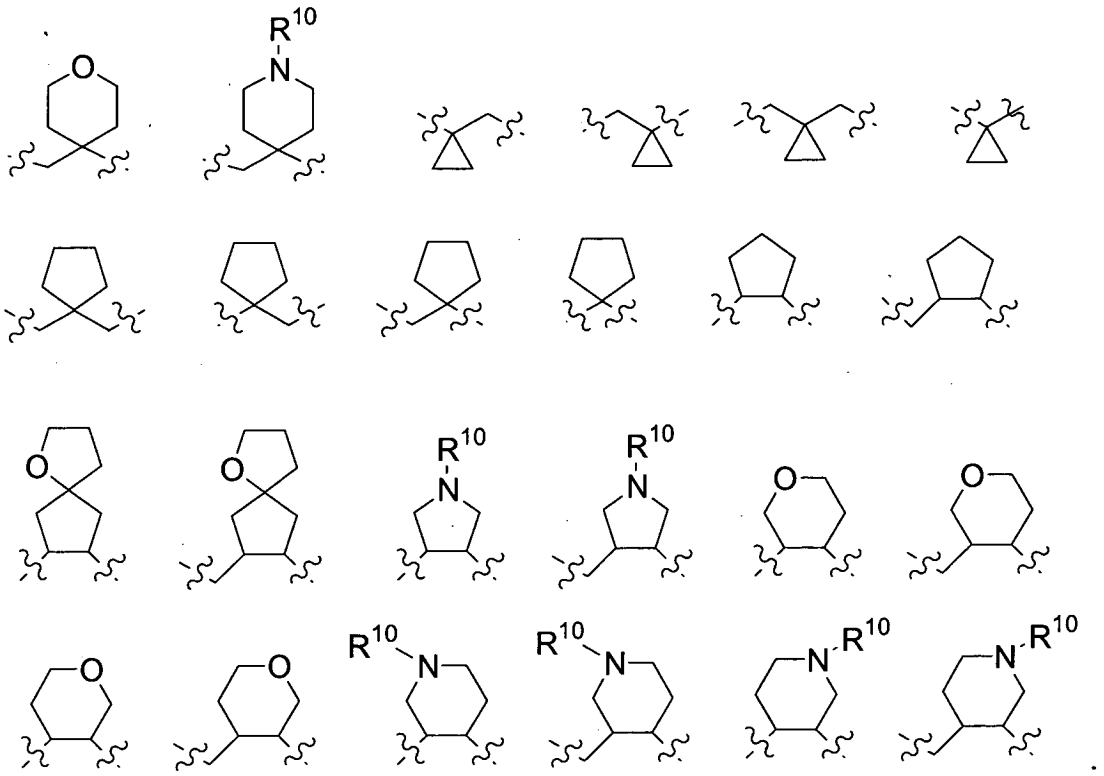
m, at each occurrence, is selected from 0, 1, and 2;

r, at each occurrence, is selected from 0, 1, and 2; and

s, at each occurrence, is selected from 0, 1, and 2.

7. (Original) A compound according to Claim 6, wherein:

L is selected from: $-CH_2-$, $-CH_2CH_2-$, $-CH_2CH_2CH_2-$, $-CH_2C(CH_3)_2-$, $-CH_2C(CH_3)_2CH_2-$,



W is $(CH_2)_m$;

Y is absent;

Z is phenyl substituted with 0-1 R^b;

Z^a is a heterocycle substituted with 0-3 R^c and selected from the group: quinolinyl, isoquinolinyl, and 1,1-dioxido-2,3-dihydro-4*H*-1,4-benzothiazin-4-yl; and

R^{10} is H, methyl, ethyl, isopropyl, isobutyl, 2-propynyl, acetyl, 2,2-dimethylpropanoyl, t-butoxycarbonyl, 3-methylbutanoyl, isobutyryl, isonicotinoyl, phenoxyacetyl, methanesulfonyl, 3-pyridinylmethyl, 4-pyridinylmethyl, 3-pyridinylcarbonyl, 4-piperidinylcarbonyl, 4-morpholinylacetyl, 4-morpholinomethyl, or [1-(t-butoxycarbonyl)-4-piperidinyl]carbonyl.

8. (Original) A compound according to Claim 1, wherein the compound is selected from the group:

4-(2-methyl-quinolin-4-ylmethoxy)-*N*-[3-(5-thioxo-4,5-dihydro-1*H*-[1,2,4]triazol-3-yl)-propyl]-benzamide;

4-(2-methyl-quinolin-4-ylmethoxy)-*N*-[2-(5-thioxo-4,5-dihydro-1*H*-[1,2,4]triazol-3-yl)-ethyl]-benzamide;

4-(2-methyl-quinolin-4-ylmethoxy)-*N*-(5-thioxo-4,5-dihydro-1*H*-[1,2,4]triazol-3-ylmethyl)-benzamide;

N-[1,1-dimethyl-2-(5-thioxo-4,5-dihydro-1*H*-[1,2,4]triazol-3-yl)-ethyl]-4-(2-methyl-quinolin-4-ylmethoxy)-benzamide;

4-[4-(2-methyl-quinolin-4-ylmethoxy)-benzoylamino]-4-(5-thioxo-4,5-dihydro-1*H*-[1,2,4]triazol-3-ylmethyl)-piperidine-1-carboxylic acid *tert*-butyl ester;

4-(2-methyl-quinolin-4-ylmethoxy)-*N*-[4-(5-thioxo-4,5-dihydro-1*H*-[1,2,4]triazol-3-ylmethyl)-tetrahydro-pyran-4-yl]-benzamide;

4-(2-methyl-quinolin-4-ylmethoxy)-*N*-[1-(5-thioxo-4,5-dihydro-1*H*-[1,2,4]triazol-3-yl)-cyclopropylmethyl]-benzamide;

5-{1-[4-(2-methyl-quinolin-4-ylmethoxy)-benzenesulfonylmethyl]-cyclopentylmethyl}-2,4-dihydro-[1,2,4]triazole-3-thione;

4-[4-(2-isopropyl-benzoimidazol-1-ylmethyl)-benzoylamino]-4-(5-thioxo-4,5-dihydro-1*H*-[1,2,4]triazol-3-ylmethyl)-piperidine-1-carboxylic acid *tert*-butyl ester;

4-(2-methyl-quinolin-4-ylmethoxy)-*N*-[*cis* 2-(5-thioxo-4,5-dihydro-1*H*-[1,2,4]triazol-3-yl)-cyclopentyl]-benzamide;

4-(2-isopropyl-benzoimidazol-1-ylmethyl)-*N*-[2-(5-thioxo-4,5-dihydro-1*H*-[1,2,4]triazol-3-yl)-cyclopentyl]-benzamide;

4-(2-methyl-quinolin-4-ylmethyl)-*N*-[2-(5-thioxo-4,5-dihydro-1*H*-[1,2,4]triazol-3-yl)-cyclopentyl]-benzamide;

4-(2-methyl-quinolin-4-ylmethoxy)-*N*-[2-(5-thioxo-4,5-dihydro-1*H*-[1,2,4]triazol-3-yl)-cyclopentyl]-benzenesulfonamide;

(3*S*,4*R*)-4-(2-methyl-quinolin-4-ylmethoxy)-*N*-[3-(5-thioxo-4,5-dihydro-1*H*-[1,2,4]triazol-3-yl)-tetrahydro-pyran-4-yl]-benzamide;

trans-4-(2-methyl-quinolin-4-ylmethoxy)-*N*-[3-(5-thioxo-4,5-dihydro-1*H*-[1,2,4]triazol-3-yl)-tetrahydro-pyran-4-yl]-benzamide;

(5*R*,7*R*,8*S*)-4-(2-methyl-quinolin-4-ylmethoxy)-*N*-[8-(5-thioxo-4,5-dihydro-1*H*-[1,2,4]triazol-3-yl)-1-oxa-spiro[4.4]non-7-yl]-benzamide;

3-[4-(2-methyl-quinolin-4-ylmethoxy)-benzoylamino]-4-(5-thioxo-4,5-dihydro-1*H*-[1,2,4]triazol-3-yl)-pyrrolidine-1-carboxylic acid *tert*-butyl ester;

4-(2-methyl-quinolin-4-ylmethoxy)-*N*-[4-(5-thioxo-4,5-dihydro-1*H*-[1,2,4]triazol-3-yl)-pyrrolidin-3-yl]-benzamide;

3-[4-(1,1-dioxo-2,3-dihydro-1*H*-1λ^6-benzo[1,4]thiazin-4-ylmethyl)-benzoylamino]-4-(5-thioxo-4,5-dihydro-1*H*-[1,2,4]triazol-3-yl)-pyrrolidine-1-carboxylic acid *tert*-butyl ester;

4-(1,1-dioxo-2,3-dihydro-1*H*-1λ^6-benzo[1,4]thiazin-4-ylmethyl)-*N*-[4-(5-thioxo-4,5-dihydro-1*H*-[1,2,4]triazol-3-yl)-pyrrolidin-3-yl]-benzamide;

(3*S*,4*S*)-4-(2-methyl-quinolin-4-ylmethoxy)-*N*-[4-(5-thioxo-4,5-dihydro-1*H*-[1,2,4]triazol-3-yl)-pyrrolidin-3-yl]-benzamide;

(3*S*,4*S*)-*N*-[1-acetyl-4-(5-thioxo-4,5-dihydro-1*H*-[1,2,4]triazol-3-yl)-pyrrolidin-3-yl]-4-(2-methyl-quinolin-4-ylmethoxy)-benzamide;

(3*S*,4*S*)-4-(2-methyl-quinolin-4-ylmethoxy)-*N*-[1-propyl-4-(5-thioxo-4,5-dihydro-1*H*-[1,2,4]triazol-3-yl)-pyrrolidin-3-yl]-benzamide;

trans-4-(2-methyl-quinolin-4-ylmethoxy)-*N*-[4-(5-thioxo-4,5-dihydro-1*H*-[1,2,4]triazol-3-yl)-pyrrolidin-3-yl]-benzamide;

(3*S*,4*S*)-3-[4-(2-methyl-quinolin-4-ylmethoxy)-benzoylamino]-4-(5-thioxo-4,5-dihydro-1*H*-[1,2,4]triazol-3-yl)-piperidine-1-carboxylic acid *tert*-butyl ester;

(3*S*,4*S*)-4-(2-methyl-quinolin-4-ylmethoxy)-*N*-[4-(5-thioxo-4,5-dihydro-1*H*-[1,2,4]triazol-3-yl)-piperidin-3-yl]-benzamide;

(3*S*,4*S*)-*N*-[1-acetyl-4-(5-thioxo-4,5-dihydro-1*H*-[1,2,4]triazol-3-yl)-piperidin-3-yl]-4-(2-methyl-quinolin-4-ylmethoxy)-benzamide;

(3*S*,4*S*)-4-(2-methyl-quinolin-4-ylmethoxy)-*N*-[1-propyl-4-(5-thioxo-4,5-dihydro-1*H*-[1,2,4]triazol-3-yl)-piperidin-3-yl]-benzamide;

(3*S*,4*S*)-*N*-[1-methanesulfonyl-4-(5-thioxo-4,5-dihydro-1*H*-[1,2,4]triazol-3-yl)-piperidin-3-yl]-4-(2-methyl-quinolin-4-ylmethoxy)-benzamide;

(3*S*,4*S*)-*N*-[1-isopropyl-4-(5-thioxo-4,5-dihydro-1*H*-[1,2,4]triazol-3-yl)-piperidin-3-yl]-4-(2-methyl-quinolin-4-ylmethoxy)-benzamide;

(3*S*,4*S*)-4-(2-methyl-quinolin-4-ylmethoxy)-*N*-[1-methyl-4-(5-thioxo-4,5-dihydro-1*H*-[1,2,4]triazol-3-yl)-piperidin-3-yl]-benzamide;

(3*S*,4*R*)-4-(2-methyl-quinolin-4-ylmethoxy)-*N*-[3-(5-thioxo-4,5-dihydro-1*H*-[1,2,4]triazol-3-yl)-piperidin-4-yl]-benzamide;

4-(2-methyl-quinolin-4-ylmethoxy)-*N*-[2-(5-oxo-4,5-dihydro-1*H*-[1,2,4]triazol-3-yl)-cyclopentyl]-benzamide;

3-[4-(2-methyl-quinolin-4-ylmethoxy)-benzoylamino]-4-(5-oxo-4,5-dihydro-1*H*-[1,2,4]triazol-3-yl)-pyrrolidine-1-carboxylic acid *tert*-butyl ester; and

4-(2-methyl-quinolin-4-ylmethoxy)-*N*-[3-(5-oxo-4,5-dihydro-1*H*-[1,2,4]triazol-3-yl)-tetrahydro-pyran-4-yl]-benzamide;

or a pharmaceutically acceptable salt form thereof.

9. (Original) A pharmaceutical composition, comprising: a pharmaceutically acceptable carrier and a therapeutically effective amount of a compound according to Claim 1 or a pharmaceutically acceptable salt form thereof.

10. (Original) A method for treating an inflammatory disorder, comprising: administering to a patient in need thereof a therapeutically effective amount of a compound according to Claim 1 or a pharmaceutically acceptable salt form thereof.

11-12. (Canceled)

13. (Original) A method of treating according to Claim 12, wherein the disease or condition is selected from to as acute infection, acute phase response, age related macular degeneration, alcoholic liver disease, allergy, allergic asthma, anorexia, aneurism, aortic aneurism, asthma, atherosclerosis, atopic dermatitis, autoimmune disease, autoimmune hepatitis, Bechet's disease, cachexia, calcium pyrophosphate dihydrate deposition disease, cardiovascular effects, chronic fatigue syndrome, chronic obstruction pulmonary disease, coagulation, congestive heart failure, corneal ulceration, Crohn's disease, enteropathic arthropathy, Felty's syndrome, fever, fibromyalgia syndrome, fibrotic disease, gingivitis, glucocorticoid withdrawal syndrome, gout, graft versus host disease, hemorrhage, HIV infection, hyperoxic alveolar injury, infectious arthritis, inflammation, intermittent hydrarthrosis, Lyme disease, meningitis, multiple sclerosis, myasthenia gravis, mycobacterial infection, neovascular

glaucoma, osteoarthritis, pelvic inflammatory disease, periodontitis, polymyositis/dermatomyositis, post-ischaemic reperfusion injury, post-radiation asthenia, psoriasis, psoriatic arthritis, pulmonary emphysema, pydoderma gangrenosum, relapsing polychondritis, Reiter's syndrome, rheumatic fever, rheumatoid arthritis, sarcoidosis, scleroderma, sepsis syndrome, Still's disease, shock, Sjogren's syndrome, skin inflammatory diseases, solid tumor growth and tumor invasion by secondary metastases, spondylitis, stroke, systemic lupus erythematosus, ulcerative colitis, uveitis, vasculitis, and Wegener's granulomatosis.

14. (Original) A method for treating inflammatory disorders, comprising: administering, to a host in need of such treatment, a therapeutically effective amount of a compound of Claim 1 or a pharmaceutically acceptable salt form thereof, in combination with one or more additional anti-inflammatory agents selected from selective COX-2 inhibitors, interleukin-1 antagonists, dihydroorotate synthase inhibitors, p38 MAP kinase inhibitors, TNF-α inhibitors and TNF-α antibody or protein sequestration agents.

15. (Original) An article of manufacture, comprising:

(a) a first container;

(b) a pharmaceutical composition located within the first container, wherein the composition, comprises: a first therapeutic agent, comprising: a compound of Claim 1 or a pharmaceutically acceptable salt form thereof; and,

(c) a package insert stating that the pharmaceutical composition can be used for the treatment of an inflammatory disorder.

16. (Original) An article of manufacture according to Claim 18, further comprising:

(d) a second container;

wherein components (a) and (b) are located within the second container and component (c) is located within or outside of the second container.

17. (Original) An article of manufacture, comprising:

(a) a first container;

(b) a pharmaceutical composition located within the first container, wherein the composition, comprises: a first therapeutic agent, comprising: a compound of Claim 1, or a pharmaceutically acceptable salt form thereof; and,

(c) a package insert stating that the pharmaceutical composition can be used in combination with a second therapeutic agent to treat an inflammatory disorder.

18. (Original) An article of manufacture according to Claim 20, further comprising:

(d) a second container;

wherein components (a) and (b) are located within the second container and component (c) is located within or outside of the second container.

19. (New) A pharmaceutical composition, comprising: a pharmaceutically acceptable carrier and a therapeutically effective amount of a compound according to Claim 2 or a pharmaceutically acceptable salt form thereof.

20. (New) A pharmaceutical composition, comprising: a pharmaceutically acceptable carrier and a therapeutically effective amount of a compound according to Claim 3 or a pharmaceutically acceptable salt form thereof.

21. (New) A pharmaceutical composition, comprising: a pharmaceutically acceptable carrier and a therapeutically effective amount of a compound according to Claim 4 or a pharmaceutically acceptable salt form thereof.

22. (New) A pharmaceutical composition, comprising: a pharmaceutically acceptable carrier and a therapeutically effective amount of a compound according to Claim 5 or a pharmaceutically acceptable salt form thereof.

23. (New) A pharmaceutical composition, comprising: a pharmaceutically acceptable carrier and a therapeutically effective amount of a compound according to Claim 6 or a pharmaceutically acceptable salt form thereof.

24. (New) A pharmaceutical composition, comprising: a pharmaceutically acceptable carrier and a therapeutically effective amount of a compound according to Claim 7 or a pharmaceutically acceptable salt form thereof.

25. (New) A pharmaceutical composition, comprising: a pharmaceutically acceptable carrier and a therapeutically effective amount of a compound according to Claim 8 or a pharmaceutically acceptable salt form thereof.